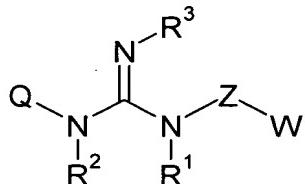


In the Claims:

1. (Previously Presented) Guanidine compound of the general formula I

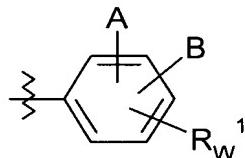


I

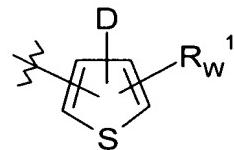
corresponding enantiomeric, diastereomeric and/or tautomeric forms thereof, as well as pharmaceutically acceptable salts thereof, wherein the given moieties have the following definitions:

W:

a moiety of the general formula W1 or W2

**W1**

or

**W2**

wherein

A:

NO_2 , NH_2 , OH , CN , CF_3 , OCF_3 , CHF_2 , OCHF_2 , COOH , $\text{O-CH}_2\text{-COOH}$,
halogen, SH , or
each optionally substituted $\text{C}_1\text{-}\text{C}_6$ -alkyl, $\text{C}_2\text{-}\text{C}_6$ -alkenyl, $\text{C}_2\text{-}\text{C}_6$ -alkynyl, $\text{C}_3\text{-}\text{C}_7$ -cycloalkyl, $\text{C}_1\text{-}\text{C}_4$ -alkylene- $\text{C}_3\text{-}\text{C}_7$ -cycloalkyl, $\text{C}_1\text{-}\text{C}_4$ -alkylene-heterocyclo-alkyl, aryl, hetaryl, heterocycloalkyl, $\text{C}_1\text{-}\text{C}_4$ -alkylene-hetaryl or $\text{C}_1\text{-}\text{C}_4$ -alkylene- aryl, or

O-R_A¹, CO-R_A¹, S-R_A¹, SO-R_A¹, CO-O-R_A¹, NR_A⁴-CO-O-R_A¹, O-CH₂-COO-R_A¹, NR_A²R_A³, CONH₂, SO₂NH₂, NR_A⁴-CO-R_A¹, SO₂-R_A¹, NR_A⁴-SO₂-R_A¹, SO₂-NR_A²R_A³ or CO-NR_A²R_A³;

R_A¹:

each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocyclo-alkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₄-alkylene-aryl, C₂-C₆-alkenylene-aryl or C₁-C₆-alkylene-hetaryl;

R_A²:

hydrogen, OH, CN, or

each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, CO-C₁-C₆-alkyl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl or SO₂-C₁-C₄-alkylene-hetaryl;

R_A³:

each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, CO-C₁-C₆-alkyl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl or SO₂-C₁-C₄-alkylene-hetaryl ;

or the the moieties $\mathbf{R_A}^2$ and $\mathbf{R_A}^3$ form, together with the nitrogen, a 3 to 7-membered, optionally substituted, saturated or aromatic heterocycle which can contain one, two or three different or same heteroatoms from the group O, N, S; wherein optionally two moieties substituted on this heterocycle can together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or same heteroatoms O, N, S, and wherein the cycle formed can optionally be substituted or a further, optionally substituted cycle can be condensed onto this cycle;

$\mathbf{R_A}^4$:

hydrogen or

each optionally substituted C₁-C₆-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₂-C₆-alkenyl, C₃-C₁₂-alkynyl, CO-C₁-C₆-alkyl, CO-O-C₁-C₆-alkyl, SO₂-C₁-C₆-alkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, CO-O-arylalkyl, CO-C₁-C₄-alkylene-aryl, CO-aryl, SO₂-aryl, hetaryl, CO-hetaryl or SO₂-C₁-C₄-alkylene-aryl;

B:

hydrogen or as moiety **A** is defined,

or each independently of one another, two of the moieties **A**, **B** or $\mathbf{R_W}^1$ form, together with a 3 to 7-membered, optionally substituted, saturated or unsaturated carbocycle or an optionally substituted, saturated or unsaturated or aromatic heterocycle which can contain one, two or three further different or same heteroatoms from the group O, N, S; wherein optionally two of the moieties substituted on this carbo- or heterocycle can together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or same heteroatoms O, N, S, and wherein the cycle formed can optionall be

substituted or a further, optionally substituted cycle can be condensed onto this cycle;

R_w¹:

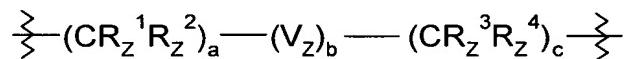
hydrogen, OH, halogen, NO₂, NH₂, CN, CF₃, CHF₂, O-CF₃, O-CHF₂, or each optionally substituted C₁-C₆-alkyl, C₃-C₇-cycloalkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₁-C₆-thioalkyl, aryl, hetaryl, O-C₁-C₆-alkyl, O-aryl, O-benzyl, C₁-C₆-alkylamino, C₁-C₆-dialkylamino, pyrrolidinyl, piperidinyl, morpholinyl, CO-C₁-C₆-alkyl, SO₂-C₁-C₆-alkyl, CO-aryl, SO₂-aryl, CO-C₁-C₄-alkylene-aryl, SO₂-C₁-C₄-alkylene-aryl, SO-aryl, CONH₂, CONH-C₁-C₆-alkyl, SO₂NH-C₁-C₆-alkyl, CON-(C₁-C₆-alkyl)₂, SO₂N-(C₁-C₆-alkyl)₂, NH-SO₂-C₁-C₆-alkyl or NH-CO-C₁-C₆-alkyl;

D:

as moiety **A** is defined;

Z:

a moiety of the general formula **Z1**



Z1

with the indices

$$a = 0 - 4$$

$$b = 0, 1$$

$$c = 0 - 4$$

wherein the sum of a, b and c is at least 1 and no more than 5;

R_Z¹, R_Z², R_Z³, R_Z⁴ independently of one another:

hydrogen, halogen, OH, or

each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₆-alkylene-C₃-C₇-cycloalkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, het-aryl or C₁-C₄-alkylene-hetaryl, or

each independently of one another, two moieties R_Z¹ and R_Z² or R_Z³ and R_Z⁴ together form a 3 to 7-membered, optionally substituted, saturated or unsaturated carbo- or heterocycle, wherein the heterocycle can contain up to three heteroatoms from the group O, N or S;

V_Z:

-CO-, -CO-NR_Z⁵-, -NR_Z⁵-CO-, -O-, -S-, -SO-, -SO₂-, -SO₂-NR_Z⁵-, -NR_Z⁵-SO₂-, -CS-, -CS-NR_Z⁵-, -NR_Z⁵-CS-, -CS-O-, -O-CS-, -CO-O-, -O-CO-, -O-, ethynylene, -C(=CR_Z⁶R_Z⁷)-, -CR_Z⁶=CR_Z⁷-, -NR_Z⁵-CO-NR_Z^{5*}-, -O-CO-NR_Z⁵-, -NR_Z⁵-;

R_Z⁵, R_Z^{5*} independently of one another:

hydrogen or

each optionally substituted C₁-C₆-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₂-C₆-alkenyl, C₃-C₁₂-alkynyl, CO-C₁-C₆-alkyl, CO-O-C₁-C₆-alkyl, SO₂-C₁-C₆-alkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-aryl, CO-aryl, SO₂-aryl, hetaryl, CO-hetaryl or SO₂-C₁-C₄-alkylene-aryl;

R_Z⁶, R_Z⁷ independently of one another:

hydrogen, OH or

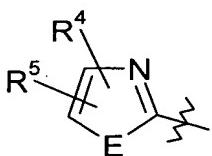
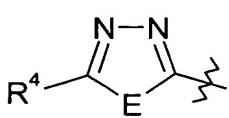
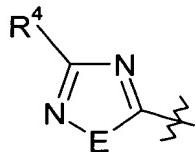
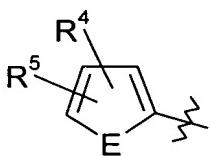
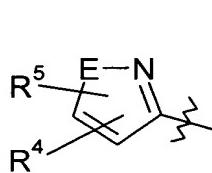
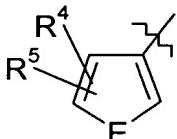
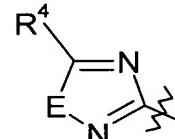
each optionally substituted C₁-C₆-alkyl, C₁-C₄-alkoxy, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₆-alkylene-C₃-C₇-cycloalkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, hetaryl or C₁-C₄-alkylene-hetaryl;

R¹, R², R³ independently of one another:

hydrogen, OH, CN, or

each optionally substituted C₁-C₆-alkyl, O-C₁-C₆-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₃-C₇-cycloalkyl, O-C₃-C₇-cycloalkyl, aryl, hetaryl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, O-aryl, O-C₁-C₄-alkylene-aryl, O-hetaryl, O-C₁-C₄-alkylene-hetaryl, CO-C₁-C₆-alkyl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-aryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl, OCO-C₁-C₆-alkyl, OCO-aryl, OCO-hetaryl, OCO-C₁-C₄-alkylene-aryl, OCO-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl or SO₂-C₁-C₄-alkylene-aryl, or

each independent from the third moiety two moieties of R¹, R² or R³ together form a 5 to 7-membered, optionally substituted, saturated or unsaturated carbocycle or an optionally substituted, saturated or unsaturated heterocycle which can contain one, two or three further different or same heteroatoms from the group O, N, S, wherein optionally two moieties substituted on this carbo- or heterocycle together can form an annellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or same heteroatoms O, N, S, and wherein the cycle formed can be optionally substituted or a further, optionally substituted cycle can be condensed onto this cycle;

Q:a doubly substituted 5-membered hetaryl moiety chosen from **Q1** to **Q7****Q1****Q2****Q3****Q4****Q5****Q6****Q7**E: O, N-R_Q¹ or S;R_Q¹:

hydrogen or

each optionally substituted C₁-C₄-alkyl, CO-C₁-C₄-alkyl, SO₂-C₁-C₄-alkyl, CO-O-C₁-C₄-alkyl, aryl, C₁-C₄-alkylene-aryl, CO-aryl, CO-hetaryl, SO₂-aryl, SO₂-hetaryl, CO-O-aryl, CO-C₁-C₄-alkylene-aryl, SO₂-C₁-C₄-alkylene-aryl or CO-O-C₁-C₄-alkylene-aryl;

R⁴, R⁵ each independently of one another a moiety chosen from the groups 1.), 2.), 3.), 4.), 5.), 6.) or 7.):

- 1.) hydrogen, halogen, CN, CF₃, CHF₂, or
each optionally substituted C₁-C₁₀-alkyl, C₂-C₁₀-alkenyl, C₂-C₁₀-alkynyl, C₃-C₇-cycloalkyl, C₁-C₆-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-

alkylene-aryl, C₁-C₄-alkylene-hetaryl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₁-C₆-alkylene-O-aryl, COO-C₁-C₄-alkyl or C₁-C₄-alkylene-COO-C₁-C₄-alkyl;

- 2.) Phenyl or naphthyl, which are each substituted with R_Q², R_Q³ and R_Q⁴, wherein

R_Q², R_Q³ and R_Q⁴ each independently of one another represent a substituent from the following group:

hydrogen, NO₂, NH₂, OH, CN, CF₃, CHF₂, OCF₃, OCHF₂, COOH, O-CH₂-COOH, SH, halogen, or
 each optionally substituted aryl, hetaryl, heterocycloalkyl, C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, C₁-C₄-alkylene-aryl or C₁-C₄-alkylene-hetaryl, or
 O-R_Q⁵, S-R_Q⁵, NR_Q⁷R_Q⁸, CO-OR_Q⁶, NR_Q⁸-CO-O-R_Q⁶, O-CH₂-COO-R_Q⁶, NR_Q⁸-CO-R_Q⁶, SO₂-R_Q⁶, NR_Q⁸-SO₂-R_Q⁶, SO₂NH₂, CONH₂, SO₂-NR_Q⁷R_Q⁸ or CO-NR_Q⁷R_Q⁸, or

two of the moieties R_Q², R_Q³ or R_Q⁴ together form a 3 to 7-membered, optionally substituted, saturated, unsaturated or aromatic carbocycle or a an optionally substituted, saturated, unsaturated aromatic heterocycle which can contain up to three further different or same heteroatoms O, N, S and optionally two of the moieties substituted on this heterocycle can together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or same heteroatoms O, N, S and the cycle formed can be optionally substituted or a further, optionally substituted cycle can be condensed onto this cycle;

- R_Q⁵** each optionally substituted C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, heterocycloalkyl or hetaryl, or C₁-C₆-alkyl, which is optionally substituted with a substituent from the group consisting of halogen, NO₂, NH₂, OH, CN, CF₃, CHF₂, OCF₃, OCHF₂, NH-(C₁-C₆-alkyl) and N(C₁-C₆-alkyl)₂;
- R_Q⁶** each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl or C₁-C₆-alkylene-O-C₁-C₆-alkyl;
- R_Q⁷** hydrogen, OH, CN, or each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, CO-C₁-C₆-alkyl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl or SO₂-C₁-C₄-alkylene-hetaryl;
- R_Q⁸** hydrogen or each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, CO-C₁-C₆-alkyl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-

O-C₁-C₆-alkyl, CO-O-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl or SO₂-C₁-C₄-alkylene-hetaryl;

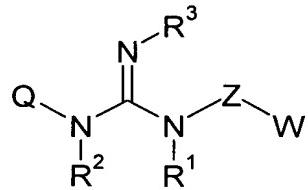
or the moieties R_Q⁷ and R_Q⁸, together with the nitrogen, form a 3 to 7-membered, optionally substituted, saturated or aromatic heterocycle, which can contain one, two or three further different or same heteroatoms O, N, S; and optionally two of the moieties substituted on this heterocycle can together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S, and the cycle formed can optionally be substituted or a further, optionally substituted cycle can be condensed onto this cycle;

- 3.) a 5- or 6-membered hetaryl moiety optionally substituted with 1 or 2 substituents, the hetaryl moiety chosen from the group consisting of:
 2-pyrrolyl, 3-pyrrolyl, 2-thiazolyl, 4-thiazolyl, 5-thiazolyl, 2-oxazolyl, 4-oxazolyl, 5-oxazolyl, 2-pyrimidyl, 4-pyrimidyl, 5-pyrimidyl, 6-pyrimidyl, 3-pyrazolyl, 4-pyrazolyl, 5-pyrazolyl, 3-isothiazolyl, 4-isothiazolyl, 5-isothiazolyl, 2-imidazolyl, 4-imidazolyl, 5-imidazolyl, 3-pyridazinyl, 4-pyridazinyl, 5-pyridazinyl, 6-pyridazinyl, 3-isoxazolyl, 4-isoxazolyl, 5-isoxazolyl, thiadiazolyl, oxadiazolyl or triazinyl or their anellated derivatives indazolyl, benzothiophenyl, benzofuranyl, indolinyl, benzimidazolyl, benzthiazolyl, benzoxazolyl, chinolinyl and isochinolinyl; or

2-thienyl or 3-thienyl optionally substituted with one or two substituents, wherein the substituents are chosen from the group consisting of halogen, NO₂, NH₂, OH, CN, CF₃, OCF₃, CHF₂, O-CHF₂, C₁-C₆-alkyl, O-C₁-C₆-alkyl, NH-(C₁-C₆-alkyl), N(C₁-C₆-alkyl)₂, NHCO-C₁-C₄-alkyl, NHSO₂-C₁-C₄-alkyl and SO₂-C₁-C₄-alkyl;

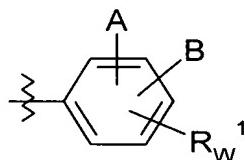
- 4.) both moieties R⁴ and R⁵ together form a 4 to 7-membered, optionally substituted, saturated or unsaturated or aromatic carbocycle or a 5- or 6-membered optionally substituted, saturated or unsaturated or aromatic heterocycle, which can contain up to three further different or identical heteroatoms O, N, S; and can be substituted with up to two further moieties, wherein optionally two moieties substituted on this carbo or hetero cycle can together form an anellated, saturated, unsaturated or aromatic carbo cycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S and wherein the cycle formed can be optionally substituted or a further, optionally substituted cycle can be condensed onto this cycle;
- 5.) a C₅-C₁₈- bi- or tricyclic, saturated hydrocarbon moiety;
- 6.) each optionally substituted C₁-C₈-Alkyl-NH₂, C₁-C₈-Alkyl-NR_Q⁷R_Q⁸, C₁-C₈-Alkyl-CO-NR_Q⁷R_Q⁸, C₁-C₈-Alkyl-SO₂NR_Q⁷R_Q⁸, C₁-C₈-Alkyl-CO-NH₂, C₁-C₈-Alkyl-SO₂NH₂, CO-NH₂, CO-NR_Q⁷R_Q⁸, SO₂NH₂, SO₂NR_Q⁷R_Q⁸, NR_Q⁷R_Q⁸;
- 7.) a 4-7-membered mono- or bicyclic saturated or unsaturated heterocycle, which can contain up to two different or identical heteroatoms from the group O, N or S, wherein this cycle can also be multiply substituted. For the case that the heterocycle contains an N-atom, this can be substituted with a moiety R_Q⁷.

2. (Previously Presented) Guanidine compound of the general formula I

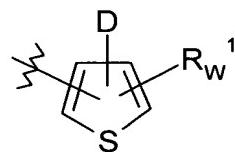


I

corresponding enantiomeric, diastereomeric and/or tautomeric forms thereof as well as pharmaceutically acceptable salts thereof, wherein the given moieties have the following definitions:

W:a moiety of the general formula **W1** or **W2**

or

**W1****W2**

wherein

A:

NO_2 , NH_2 , OH , CN , CF_3 , OCF_3 , CHF_2 , OCHF_2 , COOH , $\text{O-CH}_2\text{-COOH}$, halogen, SH , or

each optionally substituted $\text{C}_1\text{-}\text{C}_6$ -alkyl, $\text{C}_2\text{-}\text{C}_6$ -alkenyl, $\text{C}_2\text{-}\text{C}_6$ -alkynyl, $\text{C}_3\text{-}\text{C}_7$ -cycloalkyl, $\text{C}_1\text{-}\text{C}_4$ -alkylene- $\text{C}_3\text{-}\text{C}_7$ -cycloalkyl, $\text{C}_1\text{-}\text{C}_4$ -alkylene-heterocyclo-alkyl, aryl, hetaryl, heterocycloalkyl, $\text{C}_1\text{-}\text{C}_4$ -alkylene-hetaryl or $\text{C}_1\text{-}\text{C}_4$ -alkylene-aryl, or

O-R_A^1 , CO-R_A^1 , S-R_A^1 , SO-R_A^1 , CO-O-R_A^1 , $\text{NR}_\text{A}^4\text{-CO-O-R}_\text{A}^1$, $\text{O-CH}_2\text{-COO-R}_\text{A}^1$, $\text{NR}_\text{A}^2\text{R}_\text{A}^3$, CONH_2 , SO_2NH_2 , $\text{NR}_\text{A}^4\text{-CO-R}_\text{A}^1$, $\text{SO}_2\text{-R}_\text{A}^1$, $\text{NR}_\text{A}^4\text{-SO}_2\text{-R}_\text{A}^1$, $\text{SO}_2\text{-NR}_\text{A}^2\text{R}_\text{A}^3$ or $\text{CO-NR}_\text{A}^2\text{R}_\text{A}^3$;

R_A^1 :

each optionally substituted $\text{C}_1\text{-C}_6$ -alkyl, $\text{C}_2\text{-C}_6$ -alkenyl, $\text{C}_2\text{-C}_6$ -alkynyl, $\text{C}_3\text{-C}_7$ -cycloalkyl, $\text{C}_1\text{-C}_4$ -alkylene- $\text{C}_3\text{-C}_7$ -cycloalkyl, $\text{C}_1\text{-C}_4$ -alkylene-heterocyclo-alkyl, aryl, hetaryl, heterocycloalkyl, $\text{C}_1\text{-C}_4$ -alkylene-aryl, $\text{C}_2\text{-C}_6$ -alkenylene-aryl or $\text{C}_1\text{-C}_6$ -alkylene-hetaryl;

R_A^2 :

hydrogen, OH, CN, or
 each optionally substituted $\text{C}_1\text{-C}_6$ -alkyl, $\text{C}_2\text{-C}_6$ -alkenyl, $\text{C}_2\text{-C}_6$ -alkynyl, $\text{C}_1\text{-C}_4$ -alkylene- $\text{C}_3\text{-C}_7$ -cycloalkyl, $\text{C}_1\text{-C}_4$ -alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, $\text{C}_1\text{-C}_4$ -alkylene-aryl, $\text{C}_1\text{-C}_4$ -alkylene-hetaryl, $\text{C}_1\text{-C}_6$ -alkylene- $\text{O-C}_1\text{-C}_6$ -alkyl, $\text{CO-C}_1\text{-C}_6$ -alkyl, CO-aryl , CO-hetaryl , $\text{CO-C}_1\text{-C}_4$ -alkylene-aryl, $\text{CO-C}_1\text{-C}_4$ -alkylene-hetaryl, $\text{CO-O-C}_1\text{-C}_6$ -alkyl, CO-O-aryl , $\text{CO-O-C}_1\text{-C}_4$ -alkylene-aryl, CO-O-hetaryl , $\text{CO-O-C}_1\text{-C}_4$ -alkylene-hetaryl, $\text{SO}_2\text{-C}_1\text{-C}_6$ -alkyl, $\text{SO}_2\text{-aryl}$, $\text{SO}_2\text{-hetaryl}$, $\text{SO}_2\text{-C}_1\text{-C}_4$ -alkylene-aryl or $\text{SO}_2\text{-C}_1\text{-C}_4$ -alkylene-hetaryl;

R_A^3 :

each optionally substituted $\text{C}_1\text{-C}_6$ -alkyl, $\text{C}_2\text{-C}_6$ -alkenyl, $\text{C}_2\text{-C}_6$ -alkynyl, $\text{C}_1\text{-C}_4$ -alkylene- $\text{C}_3\text{-C}_7$ -cycloalkyl, $\text{C}_1\text{-C}_4$ -alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, $\text{C}_1\text{-C}_4$ -alkylene-aryl, $\text{C}_1\text{-C}_4$ -alkylene-hetaryl, $\text{C}_1\text{-C}_6$ -alkylene- $\text{O-C}_1\text{-C}_6$ -alkyl, $\text{CO-C}_1\text{-C}_6$ -alkyl, CO-aryl , CO-hetaryl , $\text{CO-C}_1\text{-C}_4$ -alkylene-aryl, $\text{CO-C}_1\text{-C}_4$ -alkylene-hetaryl, $\text{CO-O-C}_1\text{-C}_6$ -alkyl, CO-O-aryl , $\text{CO-O-C}_1\text{-C}_4$ -alkylene-aryl, CO-O-hetaryl , $\text{CO-O-C}_1\text{-C}_4$ -alkylene-hetaryl, $\text{SO}_2\text{-C}_1\text{-C}_6$ -alkyl, $\text{SO}_2\text{-aryl}$, $\text{SO}_2\text{-hetaryl}$, $\text{SO}_2\text{-C}_1\text{-C}_4$ -alkylene-aryl or $\text{SO}_2\text{-C}_1\text{-C}_4$ -alkylene-hetaryl;

or the moieties $\mathbf{R_A}^2$ and $\mathbf{R_A}^3$ form, together with the nitrogen, a 3 to 7-membered, optionally substituted, saturated or aromatic heterocycle, which can contain one, two or three further different or identical heteroatoms from the group O, N, S; wherein optionally two of the moieties substituted on this heterocycle can together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S and wherein the so-formed cycle can be optionally substituted or a further, optionally substituted cycle can be condensed onto this cycle;

$\mathbf{R_A}^4$:

hydrogen, or
each optionally substituted C₁-C₆-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₂-C₆-alkenyl, C₃-C₁₂-alkynyl, CO-C₁-C₆-alkyl, CO-O-C₁-C₆-alkyl, SO₂-C₁-C₆-alkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, CO-O-arylalkyl, CO-C₁-C₄-alkylene-aryl, CO-aryl, SO₂-aryl, hetaryl, CO-hetaryl or SO₂-C₁-C₄-alkylene-aryl;

B:

hydrogen or as moiety **A** is defined,

or each independently of one another, two of the moieties **A**, **B** or $\mathbf{R_w}^1$ together form a 3 to 7-membered, optionally substituted, saturated or unsaturated carbocycle or an optionally substituted, saturated or unsaturated or aromatic heterocycle which can contain one, two or three further different or identical heteroatoms from the group O, N, S; wherein optionally two moieties substituted on this carbo- or heterocycle can together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S and wherein the cycle formed can optionally

be substituted or a further, optionally substituted cycle can be condensed onto this cycle;

R_w¹:

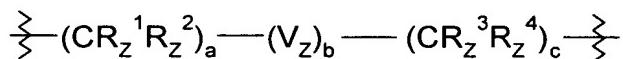
hydrogen, OH, halogen, NO₂, NH₂, CN, CF₃, CHF₂, O-CF₃, O-CHF₂, or each optionally substituted C₁-C₆-alkyl, C₃-C₇-cycloalkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₁-C₆-thioalkyl, aryl, hetaryl, O-C₁-C₆-alkyl, O-aryl, O-benzyl, C₁-C₆-alkylamino, C₁-C₆-dialkylamino, pyrrolidinyl, piperidinyl, morpholinyl, CO-C₁-C₆-alkyl, SO₂-C₁-C₆-alkyl, CO-aryl, SO₂-aryl, CO-C₁-C₄-alkylene-aryl, SO₂-C₁-C₄-alkylene-aryl, SO-aryl, CONH₂, CONH-C₁-C₆-alkyl, SO₂NH-C₁-C₆-alkyl, CON-(C₁-C₆-alkyl)₂, SO₂N-(C₁-C₆-alkyl)₂, NH-SO₂-C₁-C₆-alkyl or NH-CO-C₁-C₆-alkyl;

D:

as moiety **A** is defined;

Z:

a moiety of the general formula **Z1**



Z1

with the indices

$$a = 0 - 4$$

$$b = 0, 1$$

$$c = 0 - 4$$

wherein the sum of a, b and c is at least 1 and no more than 5;

$R_Z^1, R_Z^2, R_Z^3, R_Z^4$ independently of one another:

hydrogen, halogen, OH, or

each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₆-alkylene-C₃-C₇-cycloalkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, hetaryl or C₁-C₄-alkylene-hetaryl, or

each independently of one another two moieties are R_Z^1 and R_Z^2 or R_Z^3 and R_Z^4 together form a 3 to 7-membered, optionally substituted, saturated or unsaturated carbo- or heterocycle, wherein the heterocycle can contain up to three heteroatoms from the group O, N, or S;

V_Z :

-CO-, -CO-NR_z⁵-, -NR_z⁵-CO-, -O-, -S-, -SO-, -SO₂-, -SO₂-NR_z⁵-, -NR_z⁵-SO₂-, -CS-, -CS-NR_z⁵-, -NR_z⁵-CS-, -CS-O-, -O-CS-, -CO-O-, -O-CO-, -O-, ethynylene, -C(=CR_z⁶R_z⁷)-, -CR_z⁶=CR_z⁷-, -NR_z⁵-CO-NR_z^{5*}-, -O-CO-NR_z⁵-, -NR_z⁵-;

R_Z^5, R_Z^{5*} independently of one another:

hydrogen or

each optionally substituted C₁-C₆-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₂-C₆-alkenyl, C₃-C₁₂-alkynyl, CO-C₁-C₆-alkyl, CO-O-C₁-C₆-alkyl, SO₂-C₁-C₆-alkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-aryl, CO-aryl, SO₂-aryl, hetaryl, CO-hetaryl or SO₂-C₁-C₄-alkylene-aryl;

R_Z^6, R_Z^7 independently of one another:

hydrogen, OH or

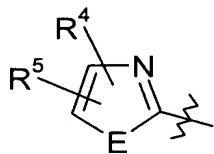
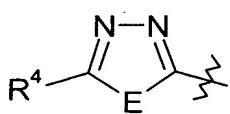
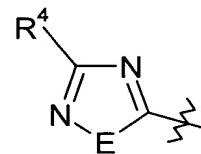
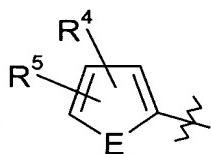
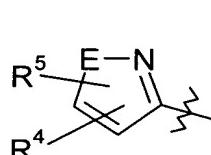
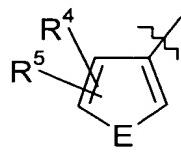
each optionally substituted C₁-C₆-alkyl, C₁-C₄-alkoxy, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₆-alkylene-C₃-C₇-cycloalkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, hetaryl or C₁-C₄-alkylene-hetaryl;

R¹, R², R³ independently of one another:

hydrogen, OH, CN, or

each optionally substituted C₁-C₆-alkyl, O-C₁-C₆-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₃-C₇-cycloalkyl, O-C₃-C₇-cycloalkyl, aryl, hetaryl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, O-aryl, O-C₁-C₄-alkylene-aryl, O-hetaryl, O-C₁-C₄-alkylene-hetaryl, CO-C₁-C₆-alkyl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-aryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl, OCO-C₁-C₆-alkyl, OCO-aryl, OCO-hetaryl, OCO-C₁-C₄-alkylene-aryl, OCO-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl or SO₂-C₁-C₄-alkylene-aryl, or

each independently of the third moiety, two moieties of **R¹, R² or R³** together form a 5 to 7-membered, optionally substituted, saturated or unsaturated carbocycle or an optionally substituted, saturated or unsaturated heterocycle which can contain one, two or three further different or identical heteroatoms from the group O, N, S, wherein optionally two of the moieties substituted on this carbo- or heterocycle can together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S and wherein the cycle formed is optionally substituted or a further, optionally substituted cycle is condensed onto this cycle;

Q:a doubly substituted 5-membered hetaryl moiety chosen from **Q1** to **Q6****Q1****Q2****Q3****Q4****Q5****Q6**E: O, N-R_Q¹ or S;R_Q¹:

hydrogen, or

each optionally substituted C₁-C₄-alkyl, CO-C₁-C₄-alkyl, SO₂-C₁-C₄-alkyl,
 CO-O-C₁-C₄-alkyl, aryl, C₁-C₄-alkylene-aryl, CO-aryl, CO-hetaryl, SO₂-
 aryl, SO₂-hetaryl, CO-O-aryl, CO-C₁-C₄-alkylene-aryl, SO₂-C₁-C₄-
 alkylene-aryl or CO-O-C₁-C₄-alkylene-aryl;

R⁴, R⁵ each independently of one another a moiety chosen from the groups 1.),
 2.), 3.), 4.) or 5.):

- 1.) hydrogen, halogen, CN, CF₃, CHF₂, or
each optionally substituted C₁-C₁₀-alkyl, C₂-C₁₀-alkenyl, C₂-C₁₀-alkynyl,
C₃-C₇-cycloalkyl, C₁-C₆-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-aryl,
C₁-C₄-alkylene-hetaryl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₁-C₆-alkylene-
O-aryl, COO-C₁-C₄-alkyl or C₁-C₄-alkylene-COO-C₁-C₄-alkyl;
- 2.) Phenyl or naphthyl, which are each substituted with R_Q², R_Q³ and R_Q⁴,
wherein

R_Q², R_Q³ and R_Q⁴ each independently of one another represent a
substituent from the following group:

hydrogen, NO₂, NH₂, OH, CN, CF₃, CHF₂, OCF₃, OCHF₂, COOH, O-
CH₂-COOH, SH, halogen, or
each optionally substituted aryl, hetaryl, heterocycloalkyl, C₁-C₆-alkyl,
C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-
cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, C₁-C₄-alkylene-aryl or C₁-
C₄-alkylene-hetaryl, or
O-R_Q⁵, S-R_Q⁵, NR_Q⁷R_Q⁸, CO-OR_Q⁶, NR_Q⁸-CO-O-R_Q⁶, O-CH₂-COO-
R_Q⁶, NR_Q⁸-CO-R_Q⁶, SO₂-R_Q⁶, NR_Q⁸-SO₂-R_Q⁶, SO₂NH₂, CONH₂, SO₂-
NR_Q⁷R_Q⁸ or CO-NR_Q⁷R_Q⁸, or

two of the moieties R_Q², R_Q³ or R_Q⁴ together form a 3 to 7-membered,
optionally substituted, saturated, unsaturated or aromatic carbocycle or
an optionally substituted, saturated or unsaturated aromatic heterocycle,
which can contain up to three further different or identical heteroatoms
O, N, S and optionally two moieties substituted on this heterocycle can
together form an anellated, saturated, unsaturated or aromatic
carbocycle or heterocycle, wherein the heterocycle can contain up to

three different or identical heteroatoms O, N, S and the cycle formed can optionally be substituted or a further, optionally substituted cycle can be condensed onto this cycle;

- R_Q⁵** each optionally substituted C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, heterocycloalkyl or hetaryl, or
C₁-C₆-alkyl, which is optionally substituted with a substituent from the group consisting of halogen, NO₂, NH₂, OH, CN, CF₃, CHF₂, OCF₃, OCHF₂, NH-(C₁-C₆-alkyl) and N(C₁-C₆-alkyl)₂;
- R_Q⁶** each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl or C₁-C₆-alkylene-O-C₁-C₆-alkyl;
- R_Q⁷** hydrogen, OH, CN, or
each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, CO-C₁-C₆-alkyl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl or SO₂-C₁-C₄-alkylene-hetaryl;
- R_Q⁸** each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, CO-C₁-C₆-alkyl, CO-aryl, CO-

hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl or SO₂-C₁-C₄-alkylene-hetaryl;

or the moieties R_Q⁷ and R_Q⁸ form, together with the nitrogen, a 3 to 7-membered, optionally substituted, saturated or aromatic heterocycle which can contain one, two or three further or different identical heteroatoms O, N, S; and optionally two moieties substituted on this heterocycle can form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S and the cycle formed can optionally be substituted or a further, optionally substituted cycle can be condensed onto this cycle;

- 3.) a 5- or 6-membered hetaryl moiety optionally substituted with one or two substituents from the group consisting of:

2-pyrrolyl, 3-pyrrolyl, 2-thiazolyl, 4-thiazolyl, 5-thiazolyl, 2-oxazolyl, 4-oxazolyl, 5-oxazolyl, 2-pyrimidyl, 4-pyrimidyl, 5-pyrimidyl, 6-pyrimidyl, 3-pyrazolyl, 4-pyrazolyl, 5-pyrazolyl, 3-isothiazolyl, 4-isothiazolyl, 5-isothiazolyl, 2-imidazolyl, 4-imidazolyl, 5-imidazolyl, 3-pyridazinyl, 4-pyridazinyl, 5-pyridazinyl, 6-pyridazinyl, 3-isoxazolyl, 4-isoxazolyl, 5-isoxazolyl, thiadiazolyl, oxadiazolyl or triazinyl or their anellated derivatives indazolyl, benzothiophenyl, benzofuranyl, indolinyl, benzimidazolyl, benzthiazolyl, benzoxazolyl, chinolinyl and isochinolinyl; or

2-thienyl or 3-thienyl optionally substituted with one or two substituents, wherein the substituents are chosen from the group consisting of halogen, NO₂, NH₂, OH, CN, CF₃, OCF₃, CHF₂, O-CHF₂, C₁-C₆-alkyl, O-C₁-C₆-alkyl, NH-(C₁-C₆-alkyl), N(C₁-C₆-alkyl)₂, NHCO-C₁-C₄-alkyl, NSO₂-C₁-C₄-alkyl and SO₂-C₁-C₄-alkyl;

[[5]] 4.) both moieties R⁴ and R⁵ together form a 4 to 7-membered, optionally substituted, saturated or unsaturated or aromatic carbocycle or a 5- or 6-membered optionally substituted, saturated or unsaturated or aromatic heterocycle, which can contain up to three further different or identical heteroatoms O, N, S and can be substituted with up to two further moieties, wherein optionally two moieties substituted on this carbo- or heterocycle can form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S and wherein the cycle formed can be optionally substituted or a further, optionally substituted cycle can be condensed onto this cycle;

[[6]] 5.) a C₅-C₁₈- bi- or tricyclic, saturated hydrocarbon moiety.

3. (Currently Amended) Guanadine compound according to claim 1 or 2, wherein the given moieties have the following definition:

W: **W1;**

A: halogen, OH, CN, CF₃, CHF₂, OCF₃, OCHF₂, or
each optionally substituted C₁-C₆-alkyl or C₂-C₆-alkenyl,
O-CH₂-COO-R_A¹, O-R_A¹, S-R_A¹, NR_A²R_A³, NR_A⁴-CO-R_A¹ or CO-
NR_A⁴R_A¹;

R_A¹: each optionally substituted C₁-C₄-alkyl, C₃-C₇-cycloalkyl, phenyl or
benzyl;

R_A²: hydrogen, or
each optionally substituted C₁-C₄-alkyl, phenyl, benzyl, phenethyl, CO-C₁-
C₄-alkyl, CO-aryl, CO-O-C₁-C₄-alkyl, SO₂-C₁-C₄-alkyl, SO₂-aryl, SO₂-
hetaryl or SO₂-C₁-C₄-alkylene-aryl;

R_A³: each optionally substituted C₁-C₄-alkyl, phenyl, benzyl, phenethyl, CO-C₁-
C₄-alkyl, CO-aryl, CO-O-C₁-C₄-alkyl, SO₂-C₁-C₄-alkyl, SO₂-aryl, SO₂-
hetaryl, or SO₂-C₁-C₄-alkylene-aryl;
or the moieties R_A² and R_A³ together form an optionally substituted 5- or
6- membered saturated or unsaturated ring, which can contain up to two
identical or different heteroatoms from the group O and N;

R_A⁴: hydrogen or an optionally substituted C₁-C₄-alkyl moiety;

B: hydrogen or as moiety A is defined;

R_W¹: hydrogen, F, Cl, CN, CF₃, O-CF₃, or
each optionally substituted C₁-C₄-alkyl, aryl, C₁-C₆-alkylamino or C₁-C₆-
dialkylamino;

in the formula **Z1** the sum of a, b and c is 1, 2 or 3;

R_Z¹, R_Z², R_Z³, R_Z⁴ independently of one another:

hydrogen, halogen, OH, optionally substituted C₁-C₆-alkyl;

V_Z: -CO-, -CO-NR_z⁵-, -NR_z⁵-CO-, -O-, -S-;

R_Z⁵: hydrogen, CH₃;

R¹, R², R³ independently of one another:

hydrogen, OH, CN, C₁-C₄-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl,
substituted aryl, benzyl, CO-C₁-C₆-alkyl, CO-aryl, CO-C₁-C₄-alkylene-
aryl, OCO-C₁-C₆-alkyl, OCO-aryl or OCO-C₁-C₄-alkylene-hetaryl;

Q is chosen from the group consisting of **Q1**, **Q2** and **Q3**;

R_Q¹: hydrogen, optionally substituted C₁-C₄-alkyl, in the aryl moiety optionally
substituted benzyl, CO-C₁-C₄-alkyl, optionally substituted benzoyl, SO₂-
C₁-C₄-alkyl or in the aryl moiety optionally substituted SO₂-aryl.

4. (Currently Amended) Guanidine compound according to ~~at least one of the claims~~ claim 1 to 3, wherein the given moieties have the following definitions:

A: OH, F, Cl, OCF₃, OCHF₂, optionally substituted C₁-C₄-alkyl, O-C₁-C₄-
alkyl or S-C₁-C₄-alkyl;

B: hydrogen, OH, F, Cl, CF₃, OCF₃, OCHF₂, optionally substituted C₁-C₄-
alkyl, O-C₁-C₄-alkyl or S-C₁-C₄-alkyl;

R_w¹: hydrogen, F, Cl, CN, CF₃ or O-CF₃;

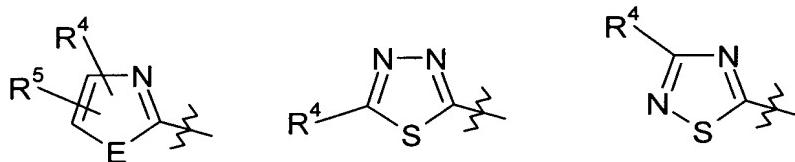
Z: each optionally substituted C₁-C₄-alkyl or C₁-C₄-alkylene-O-C₁-C₄- alkyl;

R_Z¹, R_Z², R_Z³, R_Z⁴ each independently of one another:
hydrogen, F, CH₃;

R¹, R², R³ independently of one another:

hydrogen, OH, CN, O-methyl, O-phenyl, acetyl, benzoyl, O-acetyl, O-benzoyl;

Q is chosen from the group consisting of



R_Q¹: hydrogen, CH₃, methanesulfonyl, phenylsulfonyl or tosyl.

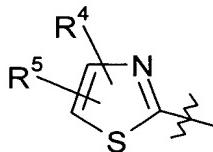
5. (Currently Amended) Guanidine compound according to at least one of the claims claim 1 to 4, wherein the given moieties have the following definitions:

A: OH, OCF₃, OCH₃, O-ethyl, O-propyl or O-i-propyl;

Z: -CH₂-, -CH₂-O-, -CH₂-CH₂- or -CH₂-CH₂-O-;

two of the moieties \mathbf{R}^1 , \mathbf{R}^2 , or \mathbf{R}^3 , are hydrogen, and the third moiety is hydrogen, OH, acetyl or benzoyl;

Q:



6. (Currently Amended) Guanidine compound according to ~~at least one of the claims~~ claim 1 to 5, wherein \mathbf{R}^4 and/or \mathbf{R}^5 each independently of one another represents a moiety chosen from the groups 1.), 2.), 3.), 4.) or 5.):

- 1.) hydrogen, F, Cl, CN, CF₃, or
each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl or C₃-C₇-cycloalkyl;
- 2.) \mathbf{R}_Q^1 , \mathbf{R}_Q^2 and \mathbf{R}_Q^3 independently of one another
hydrogen, CN, CF₃, CHF₂, OCF₃, OCHF₂, F, Cl, OH or
each optionally substituted phenyl or hetaryl, C₁-C₄-alkyl, C₅-C₇-cycloalkyl, O-R_Q⁵, NR_Q⁷R_Q⁸, CO-OR_Q⁶, NR_Q⁸-CO-O-R_Q⁶, O-CH₂-COO-R_Q⁶, NR_Q⁸-CO-R_Q⁶, SO₂-R_Q⁶, NR_Q⁸-SO₂-R_Q⁶, NR_Q⁸-CO-O-R_Q⁶, SO₂NH₂, CONH₂, SO₂-NR_Q⁷R_Q⁸ or CO-NR_Q⁷R_Q⁸;

\mathbf{R}_Q^5 : C₁-C₄-Alkyl, which is optionally substituted with a substituent from the group consisting of F, Cl, OH, CN, CF₃, OCF₃, NH-(C₁-C₄-alkyl) and N(C₁-C₄-alkyl)₂;

\mathbf{R}_Q^6 : each optionally substituted C₁-C₆-alkyl, aryl, hetaryl or phenyl;

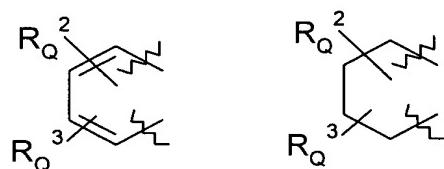
R_Q⁷: hydrogen, each optionally substituted C₁-C₄-alkyl, allyl, aryl, hetaryl, benzyl, phenethyl or CH₂-hetaryl;

R_Q⁸: each optionally substituted C₁-C₄-alkyl, allyl, aryl, hetaryl, benzyl, phenethyl or CH₂-hetaryl;

or R_Q⁷ und R_Q⁸ form an optionally substituted 3- or 7-membered saturated or unsaturated ring which can contain up to two identical or differnt hetero atoms from the group O and N;

3.) benzothiophenyl, benzofuranyl, chinolinyl or isochinolinyl;

4.) both moieties R⁴ and R⁵ together form one of the following rings:



wherein R_Q² and R_Q³ are as defined under 2.);

5.) Adamantyl.

7. (Currently Amended) Guanidine compound according to claim 1 or 2, wherein the given moieties have the following definitions:

W: W1;

A: halogen, OH, CN, CF₃, CHF₂, OCF₃, OCHF₂, or

each optionally substituted C₁-C₆-alkyl or C₂-C₆-alkenyl,
O-CH₂-COO-R_A¹, O-R_A¹, S-R_A¹, NR_A²R_A³, NR_A⁴-CO-R_A¹, SO₂NH₂,
NR_A⁴-SO₂-R_A¹, SO₂-NR_A²R_A³ or CO-NR_A⁴R_A¹;

R_A¹: each optionally substituted C₁-C₄-alkyl, C₃-C₇-cycloalkyl, phenyl or benzyl;

R_A²: hydrogen, or
each optionally substituted C₁-C₄-alkyl, phenyl, benzyl, phenethyl, CO-C₁-C₄-alkyl, CO-aryl, CO-O-C₁-C₄-alkyl, SO₂-C₁-C₄-alkyl, SO₂-aryl, SO₂-hetaryl or SO₂-C₁-C₄-alkylene-aryl;

R_A³: each optionally substituted C₁-C₄-alkyl, phenyl, benzyl, phenethyl, CO-C₁-C₄-alkyl, CO-aryl, CO-O-C₁-C₄-alkyl, SO₂-C₁-C₄-alkyl, SO₂-aryl, SO₂-hetaryl, or SO₂-C₁-C₄-alkylene-aryl;
or the moieties **R_A**² and **R_A**³ together form an optionally substituted 5- or 6-membered saturated or unsaturated ring, which can contain up to two identical or different heteroatoms from the group O and N;

R_A⁴: hydrogen or an optionally substituted C₁-C₄-alkyl moiety;

B: hydrogen or as moiety **A** is defined;

R_w¹: hydrogen, F, Cl, CN, CF₃, O-CF₃, or
each optionally substituted C₁-C₄-alkyl, aryl, C₁-C₆-alkylamino or C₁-C₆-dialkylamino;

in the formula **Z1** the sum of a, b and c is 1, 2 or 3;

R_Z¹, **R_Z**², **R_Z**³, **R_Z**⁴ independently of one another:

hydrogen, halogen, OH, optionally substituted C₁-C₆-alkyl;

V_Z: -CO-, -CO-NR_z⁵-, -NR_z⁵-CO-, -O-, -S-;

R_Z⁵: hydrogen, CH₃;

R¹, R², R³ independently of one another:

hydrogen, OH, CN, C₁-C₄-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl,
substituted aryl, benzyl, CO-C₁-C₆-alkyl, CO-aryl, CO-C₁-C₄-alkylene-
aryl, OCO-C₁-C₆-alkyl, OCO-aryl or OCO-C₁-C₄-alkylene-hetaryl;

Q is chosen from the group consisting of **Q1, Q2, Q3** and **Q5**;

R_Q¹: hydrogen, optionally substituted C₁-C₄-alkyl, in the aryl moiety optionally
substituted benzyl, CO-C₁-C₄-alkyl, optionally substituted benzoyl, SO₂-
C₁-C₄-alkyl or in the aryl moiety optionally substituted SO₂-aryl.

8. (Currently Amended) Guanidine compound according to ~~at least one of claims~~
~~claim 1, 2 or 7~~, wherein the given moieties have the following definitions:

A: OH, F, Cl, OCF₃, OCHF₂, optionally substituted C₁-C₄-alkyl, O-C₁-C₄-
alkyl or S-C₁-C₄-alkyl;

B: hydrogen, OH, F, Cl, CF₃, OCF₃, OCHF₂, optionally substituted C₁-C₄-
alkyl, O-C₁-C₄-alkyl or S-C₁-C₄-alkyl;

R_W¹: hydrogen, F, Cl, CN, CF₃ or O-CF₃;

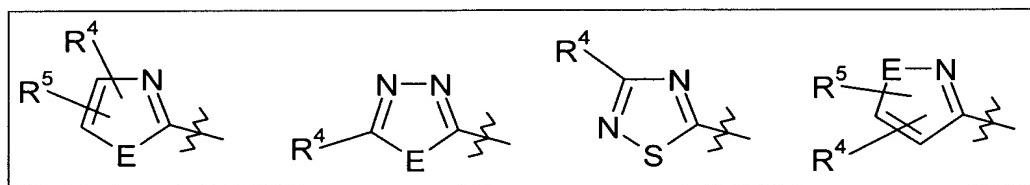
Z: each optionally substituted C₁-C₄-alkyl or C₁-C₄-alkylene-O-C₁-C₄-alkyl;

R_Z¹, R_Z², R_Z³, R_Z⁴ each independently of one another:
hydrogen, F, CH₃;

R¹, R², R³ independently of one another:

hydrogen, OH, CN, O-methyl, O-phenyl, acetyl, benzoyl, O-acetyl, O-benzoyl;

Q is chosen from the group consisting of



R_Q¹: hydrogen, CH₃, phenyl, benzyl, methanesulfonyl, phenylsulfonyl or tosyl.

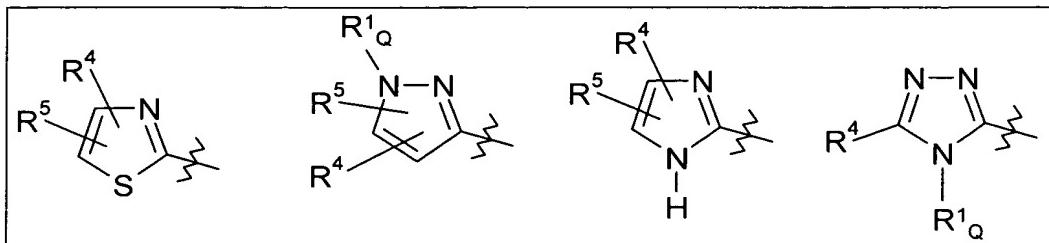
9. (Currently Amended) Guanidine compound according to ~~at least one of the claims claim 1, 2, 7 or 8~~, wherein the given moieties have the following definitions:

A: OH, OCF₃, OCH₃, O-ethyl, O-propyl or O-i-propyl;

Z: -CH₂-, -CH₂-O-, -CH₂-CH₂- or -CH₂-CH₂-O-;

two of the moieties **R¹, R², or R³** are hydrogen, and the third moiety is hydrogen, OH, acetyl or benzoyl;

Q:



\mathbf{R}_Q^1 : hydrogen, CH₃, phenyl, benzyl, methanesulfonyl, phenylsulfonyl or tosyl.

10. (Currently Amended) Guanidine compound according to ~~at least one of the claims claim 1, 2 or 7 to 9~~, wherein \mathbf{R}^4 and/or \mathbf{R}^5 each independently from one another represent a moiety chosen from the groups 1.), 2.), 3.), 4.), 5.) or 7.):

- 1.) hydrogen, F, Cl, CN, CF₃, or
each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl or C₃-C₇-cycloalkyl;
- 2.) \mathbf{R}_Q^1 , \mathbf{R}_Q^2 and \mathbf{R}_Q^3 independently of one another
hydrogen, CN, CF₃, CHF₂, OCF₃, OCHF₂, F, Cl, OH or
each optionally substituted phenyl or hetaryl, C₁-C₄-alkyl, C₅-C₇-cycloalkyl, O-R_Q⁵, NR_Q⁷R_Q⁸, CO-OR_Q⁶, NR_Q⁸-CO-O-R_Q⁶, O-CH₂-COO-R_Q⁶, NR_Q⁸-CO-R_Q⁶, SO₂-R_Q⁶, NR_Q⁸-SO₂-R_Q⁶, NR_Q⁸-CO-O-R_Q⁶, SO₂NH₂, CONH₂, SO₂-NR_Q⁷R_Q⁸ or CO-NR_Q⁷R_Q⁸;

\mathbf{R}_Q^5 : C₁-C₄-alkyl, which is optionally substituted with a substituent from the group consisting of F, Cl, OH, CN, CF₃, OCF₃, NH-(C₁-C₄-alkyl) and N(C₁-C₄-alkyl)₂;

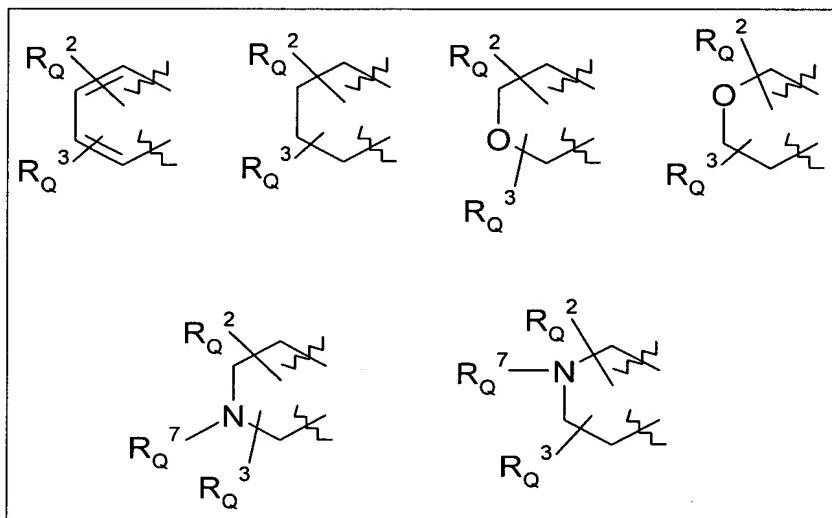
\mathbf{R}_Q^6 : each optionally substituted C₁-C₆-alkyl, aryl, hetaryl or phenyl;

R_Q^7 : hydrogen, each optionally substituted C₁-C₄-alkyl, allyl, aryl, hetaryl, benzyl, phenethyl or CH₂-hetaryl;

R_Q^8 : hydrogen, each optionally substituted C₁-C₄-alkyl, allyl, aryl, hetaryl, benzyl, phenethyl or CH₂-hetaryl;

or R_Q^7 and R_Q^8 form an optionally substituted 3- or 7-membered saturated or unsaturated ring, which can contain up to two identical or different heteroatoms from the group O and N;

- 3.) benzothiophenyl, benzofuranyl, chinolinyl or isochinolinyl;
- 4.) both moieties R^4 and R^5 together form one of the following rings:

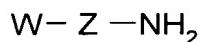


wherein R_Q^2 and R_Q^3 are defined as under 2.); or together can form an anellated 5- or 6-membered ring;

- 5.) adamantyl;

[[7]] **6.)** each optionally substituted azetidine-3-yl, pyrrolidine-2-yl, pyrrolidine-3-yl, piperidine-2-yl, piperidine-3-yl, piperidine-4-yl, tetrahydro-2H-pyrane-4-yl, tetrahydrofuran-3-yl, azepan-4-yl, azepan-3-yl, azepan-2-yl, 1,4-diazepane-5-yl, 1,2,3,6-tetrahydropyridine-4-yl, 2,5-dihydro-1H-pyrrol-3-yl.

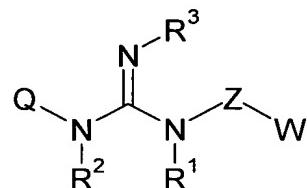
11. (Currently Amended) Guanidine compound according to ~~at least one of the claims claim 1-to-10~~, wherein one moiety from \mathbf{R}^4 and \mathbf{R}^5 is chosen from group 1.), and the other moiety from \mathbf{R}^4 and \mathbf{R}^5 is chosen from the group 1.), 2.) or 3.).
12. (Cancel)
13. (Currently Amended) Pharmaceutical composition, comprising at least one guanidine compound according to ~~one of the claims claim 1-to-12~~, as well as a pharmaceutically acceptable carrier or dilution agent.
14. (Currently Amended) ~~Use of compounds of the general formula IVA A method for the preparation of 5HT5A receptor ligands comprising using a compound of the general formula IVA:~~



IVA

15. (Currently Amended) ~~Use A method according to claim 14 wherein the 5HT5A receptor ligand is a for the preparation of the guanidine compounds compound according to one of the claims claim 1-to-12.~~

16. (Currently Amended) Use A method of treating a patient having a disease modulated by 5-HT₅ receptor activity comprising administering to said patient an effective amount of a guanidine compound of the general formula IA



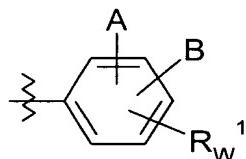
IA

of the corresponding enantiomeric, diastereomeric and/or tautomeric forms thereof as well as pharmaceutically acceptable salts thereof

wherein the given moieties have the following definitions:

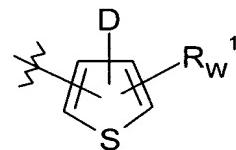
W:

a moiety of the general formula W1 or W2



W1

or



W2

A:

NO₂, NH₂, OH, CN, CF₃, OCF₃, CHF₂, OCHF₂, COOH, O-CH₂-COOH, halogen, SH, or

each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl or C₁-C₄-alkylene-hetero-cycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₄-alkylene-hetaryl or C₁-C₄-alkylene-aryl, or

O-R_A¹, CO-R_A¹, S-R_A¹, SO-R_A¹, CO-O-R_A¹, NR_A⁴-CO-O-R_A¹, O-CH₂-COO-R_A¹, NR_A²R_A³, CONH₂, SO₂NH₂, NR_A⁴-CO-R_A¹, SO₂-R_A¹, NR_A⁴-SO₂-R_A¹, SO₂-NR_A²R_A³ or CO-NR_A²R_A³;

R_A¹:

each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocyclo-alkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₄-alkylene-aryl, C₂-C₆-alkenylene-aryl or C₁-C₆-alkylene-hetaryl;

R_A²:

hydrogen, OH, CN, or
 each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, CO-C₁-C₆-alkyl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl or SO₂-C₁-C₄-alkylene-hetaryl;

R_A³:

each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, CO-C₁-C₆-alkyl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl or SO₂-C₁-C₄-alkylene-hetaryl;

or the moieties \mathbf{R}_A^2 and \mathbf{R}_A^3 form, together with the nitrogen, a 3 to 7-membered, optionally substituted, saturated or aromatic heterocycle, which can contain one, two or three further different or identical heteroatoms from the group O, N, S; wherein optionally two moieties substituted on this heterocycle can together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms, O, N, S and wherein the so-formed cycle can optionally be substituted or a further, optionally substituted cycle can be condensed onto this cycle;

\mathbf{R}_A^4 :

hydrogen, or

each optionally substituted C₁-C₆-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₂-C₆-alkenyl, C₃-C₁₂-alkynyl, CO-C₁-C₆-alkyl, CO-O-C₁-C₆-alkyl, SO₂-C₁-C₆-alkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, CO-O-arylalkyl, CO-C₁-C₄-alkylene-aryl, CO-aryl, SO₂-aryl, hetaryl, CO-hetaryl or SO₂-C₁-C₄-alkylene-aryl;

B:

hydrogen or as moiety **A** is defined,

or each independently of one another two of the moieties **A**, **B** or \mathbf{R}_W^1 together form a 3 to 7-membered, optionally substituted, saturated or unsaturated carbocycle with an optionally substituted, saturated or unsaturated or aromatic heterocycle which can contain one, two or three further different or identical heteroatoms from the group O, N, S; wherein optionally two moieties substituted on this carbo- or heterocycle can form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S and wherein the cycle formed can optionally be

substituted or a further, optionally substituted cycle can be condensed onto this cycle;

R_W¹:

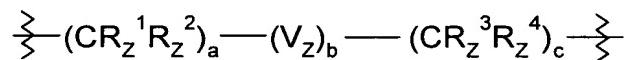
hydrogen, OH, halogen, NO₂, NH₂, CN, CF₃, CHF₂, O-CF₃, O-CHF₂, or each optionally substituted C₁-C₆-alkyl, C₃-C₇-cycloalkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₁-C₆-alkylene-S-C₁-C₆-alkyl, aryl, hetaryl, O-C₁-C₆-alkyl, O-aryl, O-benzyl, C₁-C₆-alkylamino, C₁-C₆-dialkylamino, pyrrolidinyl, piperidinyl, morpholinyl, CO-C₁-C₆-alkyl, SO₂-C₁-C₆-alkyl, CO-aryl, SO₂-aryl, CO-C₁-C₄-alkylene-aryl, SO₂-C₁-C₄-alkylene-aryl, SO-aryl, CONH₂, CONH-C₁-C₆-alkyl, SO₂NH-C₁-C₆-alkyl, CON(C₁-C₆-alkyl)₂, SO₂N(C₁-C₆-alkyl)₂, NH-SO₂-C₁-C₆-alkyl or NH-CO-C₁-C₆-alkyl;

D:

as moiety **A** is defined;

Z:

a moiety of the general formula **Z1**



Z1

with the indices

$$a = 0 - 4$$

$$b = 0, 1$$

$$c = 0 - 4$$

wherein the sum of a, b and c is no more than 5;

R_Z¹, R_Z², R_Z³, R_Z⁴ independently of one another:

hydrogen, halogen, OH, or

each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₆-alkylene-C₃-C₇-cycloalkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, hetaryl or C₁-C₄-alkylene-hetaryl, or

each independently of one another two moieties R_Z¹ and R_Z² or R_Z³ and R_Z⁴ together form a 3 to 7-membered, optionally substituted, saturated or unsaturated carbo- or heterocycle, which can contain up to three heteroatoms from the group O, N or S;

V_Z:

-CO-, -CO-NR_z⁵-, -NR_z⁵-CO-, -O-, -S-, -SO-, -SO₂-, -SO₂-NR_z⁵-, -NR_z⁵-SO₂-, -CS-, -CS-NR_z⁵-, -NR_z⁵-CS-, -CS-O-, -O-CS-, -CO-O-, -O-CO-, -O-, ethynylene, -C(=CR_Z⁶R_z⁷)-, -CR_Z⁶=CR_Z⁷-, -NR_z⁵-CO-NR_z^{5*}-, -O-CO-NR_z⁵-, -NR_z⁵-;

R_Z⁵, R_Z^{5*} independently of one another:

hydrogen, or

each optionally substituted C₁-C₆-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₂-C₆-alkenyl, C₃-C₁₂-alkynyl, CO-C₁-C₆-alkyl, CO-O-C₁-C₆-alkyl, SO₂-C₁-C₆-alkyl, C₃-C₇-cycloalkyl, Aryl, C₁-C₄-alkylene-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-aryl, CO-aryl, SO₂-aryl, hetaryl, CO-hetaryl or SO₂-C₁-C₄-alkylene-aryl;

R_Z⁶, R_Z⁷ independently of one another:

hydrogen, OH, or

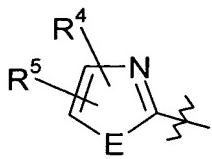
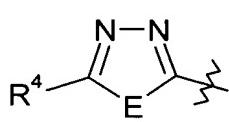
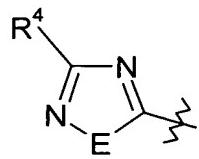
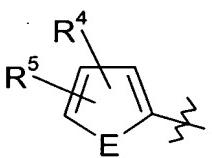
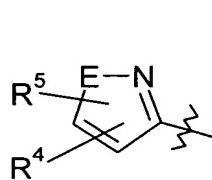
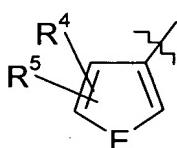
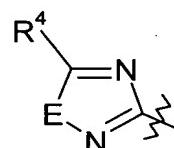
each optionally substituted C₁-C₆-alkyl, C₁-C₄-alkoxy, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₆-alkylene-C₃-C₇-cycloalkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, hetaryl or C₁-C₄-alkylene-hetaryl;

R¹, R², R³ independently of one another:

hydrogen, OH, CN, or

each optionally substituted C₁-C₆-alkyl, O-C₁-C₆-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₃-C₇-cycloalkyl, O-C₃-C₇-cycloalkyl, aryl, hetaryl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, O-aryl, O-C₁-C₄-alkylene-aryl, O-hetaryl, O-C₁-C₄-alkylene-hetaryl, CO-C₁-C₆-alkyl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-aryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-Aryl, OCO-C₁-C₆-alkyl, OCO-aryl, OCO-hetaryl, OCO-C₁-C₄-alkylene-aryl, OCO-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl or SO₂-C₁-C₄-alkylene-aryl, or

each independently from the third moiety two moieties of R¹, R² or R³ together form a 5 to 7-membered, optionally substituted, saturated or unsaturated carbocycle or an optionally substituted, saturated or unsaturated, which can contain one, two or three further different or identical heteroatoms from the group O, N, S, wherein optionally two moieties substituted on this carbo- or heterocycle can together form an annellated, saturated, unsaturated, or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S and wherein the cycle formed can be optionally substituted or a further, optionally substituted cycle can be condensed onto this cycle;

Q:a doubly substituted 5-membered hetaryl moiety, chosen from **Q1** to **Q7****Q1****Q2****Q3****Q4****Q5****Q6****Q7**

E: O, N-R_Q¹ or S;

R_Q¹:

hydrogen, or

each optionally substituted C₁-C₄-alkyl, CO-C₁-C₄-alkyl, SO₂-C₁-C₄-alkyl,
CO-O-C₁-C₄-alkyl, aryl, C₁-C₄-alkylene-aryl, CO-aryl, CO-hetaryl, SO₂-
aryl, SO₂-hetaryl, CO-O-aryl, CO-C₁-C₄-alkylene-aryl, SO₂-C₁-C₄-
alkylene-aryl or CO-O-C₁-C₄-alkylene-aryl;

R⁴, R⁵ each independently of one another, a moiety chosen from the groups 1.),
2.), 3.), 4.), 5.), 6.) or 7.):

1.) hydrogen, halogen, CN, CF₃, CHF₂, or
 each optionally substituted C₁-C₁₀-alkyl, C₂-C₁₀-alkenyl, C₂-C₁₀-alkynyl, C₃-C₇-cycloalkyl, C₁-C₆-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₁-C₆-alkylene-O-aryl, COO-C₁-C₄-alkyl or C₁-C₄-alkylene-COO-C₁-C₄-alkyl;

[[3]] 2.) phenyl or naphthyl, which are substituted with R_Q², R_Q³ and R_Q⁴,
 wherein

R_Q², R_Q³ and R_Q⁴ each independently of one another represent a
 substituent from the following group:

hydrogen, NO₂, NH₂, OH, CN, CF₃, CHF₂, OCF₃, OCHF₂, COOH, O-CH₂-COOH, SH, halogen, or
 each optionally substituted aryl, hetaryl, heterocycloalkyl, C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, C₁-C₄-alkylene-aryl or C₁-C₄-alkylene-hetaryl, or
 O-R_Q⁵, S-R_Q⁵, NR_Q⁷R_Q⁸, CO-OR_Q⁶, NR_Q⁸-CO-O-R_Q⁶, O-CH₂-COO-R_Q⁶, NR_Q⁸-CO-R_Q⁶, SO₂-R_Q⁶, NR_Q⁸-SO₂-R_Q⁶, SO₂NH₂, CONH₂, SO₂-NR_Q⁷R_Q⁸ or CO-NR_Q⁷R_Q⁸, or

two of the moieties R_Q², R_Q³ or R_Q⁴ together form a 3 to 7-membered, optionally substituted, saturated, unsaturated carbocycle or an optionally substituted, saturated, unsaturated aromatic heterocycle which can contain up to three further different or identical heteroatoms O, N, S, and optionally two moieties substituted on this heterocycle can form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S, and the cycle formed may optionally be

substituted or a further, optionally substituted cycle can be condensed onto this cycle;

R_Q⁵ each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, heterocycloalkyl, aryl or hetaryl;

R_Q⁶ each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl or C₁-C₆-alkylene-O-C₁-C₆-alkyl;

R_Q⁷ hydrogen, OH, CN, or
each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, CO-C₁-C₆-alkyl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl or SO₂-C₁-C₄-alkylene-hetaryl;

R_Q⁸ hydrogen or
each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-

C₆-alkylene-O-C₁-C₆-alkyl, CO-C₁-C₆-alkyl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl or SO₂-C₁-C₄-alkylene-hetaryl;

or both moieties R_Q⁷ and R_Q⁸ form, together with the nitrogen, a 3 to 7-membered, optionally substituted, saturated or aromatic heterocycle, which can contain one, two or three different or identical heteroatoms O, N, S; and optionally two moieties substituted on this heterocycle can together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S, and the cycle formed can be optionally substituted or a further, optionally substituted cycle can be condensed onto this cycle;

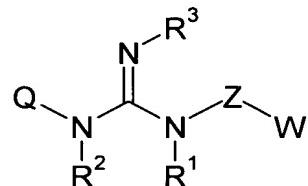
- 3.) a 5- or 6-membered, hetaryl moiety, optionally substituted with 1 or 2 substituents from the group consisting of:

2-furyl, 3-furyl, 2-pyrrolyl, 3-pyrrolyl, 2-thienyl, 3-thienyl, 2-pyridyl, 3-pyridyl, 4-pyridyl, 2-thiazolyl, 4-thiazolyl, 5-thiazolyl, 2-oxazolyl, 4-oxazolyl, 5-oxazolyl, 2-pyrimidyl, 4-pyrimidyl, 5-pyrimidyl, 6-pyrimidyl, 3-pyrazolyl, 4-pyrazolyl, 5-pyrazolyl, 3-isothiazolyl, 4-isothiazolyl, 5-isothiazolyl, 2-imidazolyl, 4-imidazolyl, 5-imidazolyl, 3-pyridazinyl, 4-pyridazinyl, 5-pyridazinyl, 6-pyridazinyl, 3-isoxazolyl, 4-isoxazolyl, 5-isoxazolyl, thiadiazolyl, oxadiazolyl or triazinyl or their anellated derivatives indazolyl, indolyl, benzothiophenyl, benzofuranyl, indolinyl, benzimidazolyl, benzthiazolyl, benzoxazolyl, chinolinyl and isochinolinyl;

- 4.) both moieties \mathbf{R}^4 and \mathbf{R}^5 together form a 4 to 7-membered, optionally substituted, saturated or unsaturated or aromatic carbocycle or a 5- or 6-membered optionally substituted, saturated or unsaturated or aromatic heterocycle, which can contain up to three further different or identical heteroatoms O, N, S, and which can be substituted with up to two further moieties, wherein optionally two moieties substituted on this carbo- or heterocycle together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S and wherein the cycle formed can optionally be substituted or a further, optionally substituted cycle can be condensed onto this cycle;
- 5.) a C₆-C₁₀- bi- or tricyclic, saturated hydrocarbon moiety;
- 6.) each optionally substituted C₁-C₈-alkyl-NH₂, C₁-C₈-alkyl-NR_Q⁷R_Q⁸, C₁-C₈-alkyl-CO-NR_Q⁷R_Q⁸, C₁-C₈-alkyl-SO₂NR_Q⁷R_Q⁸, C₁-C₈-alkyl-CO-NH₂, C₁-C₈-alkyl-SO₂NH₂, CO-NH₂, CO-NR_Q⁷R_Q⁸, SO₂NH₂, SO₂NR_Q⁷R_Q⁸, NR_Q⁷R_Q⁸;
- 7.) a 4-7-membered mono- or bicyclic saturated or unsaturated heterocycle, which can contain up to two different or identical heteroatoms from the group O, N or S, wherein this cycle can also be multiply substituted. For the case that the heterocycle contains an N-atom, this can be substituted with an R_Q⁷ moiety.

~~for the preparation of a medicament for the treatment of diseases modulated by a 5-HT₅ receptor activity.~~

17. (Currently Amended) Use A method of treating a patient having a disease modulated by 5-HT₅ receptor activity comprising administering to said patient and effective amount of a guanidine compound of the general formula IA



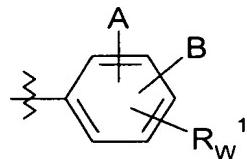
IA

of the corresponding enantiomeric, diastereomeric and/or tautomeric forms thereof as well as pharmaceutically acceptable salts thereof,

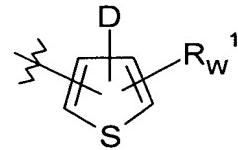
wherein the given moieties have the following definitions:

W:

a moiety of the general formula W1 or W2



or



W2

A:

NO₂, NH₂, OH, CN, CF₃, OCF₃, CHF₂, OCHF₂, COOH, O-CH₂-COOH, halogen, SH, or
each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl or C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₄-alkylene-hetaryl or C₁-C₄-alkylene-aryl, or

$O-R_A^1$, $CO-R_A^1$, $S-R_A^1$, $SO-R_A^1$, $CO-O-R_A^1$, $NR_A^4-CO-O-R_A^1$, $O-CH_2-COO-R_A^1$, $NR_A^2R_A^3$, $CONH_2$, SO_2NH_2 , $NR_A^4-CO-R_A^1$, $SO_2-R_A^1$, $NR_A^4-SO_2-R_A^1$, $SO_2-NR_A^2R_A^3$ or $CO-NR_A^2R_A^3$;

R_A^1 :

each optionally substituted C_1-C_6 -alkyl, C_2-C_6 -alkenyl, C_2-C_6 -alkynyl, C_3-C_7 -cycloalkyl, C_1-C_4 -alkylene- C_3-C_7 -cycloalkyl, C_1-C_4 -alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C_1-C_4 -alkylene-aryl, C_2-C_6 -alkenylene-aryl or C_1-C_6 -alkylene-hetaryl;

R_A^2 :

hydrogen, OH, CN, or

each optionally substituted C_1-C_6 -alkyl, C_2-C_6 -alkenyl, C_2-C_6 -alkynyl, C_1-C_4 -alkylene- C_3-C_7 -cycloalkyl, C_1-C_4 -alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C_1-C_4 -alkylene-aryl, C_1-C_4 -alkylene-hetaryl, C_1-C_6 -alkylene-O- C_1-C_6 -alkyl, $CO-C_1-C_6$ -alkyl, CO -aryl, CO -hetaryl, $CO-C_1-C_4$ -alkylene-aryl, $CO-C_1-C_4$ -alkylene-hetaryl, $CO-O-C_1-C_6$ -alkyl, CO -aryl, $CO-O-C_1-C_4$ -alkylene-aryl, $CO-O$ -hetaryl, $CO-O-C_1-C_4$ -alkylene-hetaryl, $SO_2-C_1-C_6$ -alkyl, SO_2 -aryl, SO_2 -hetaryl, $SO_2-C_1-C_4$ -alkylene-aryl or $SO_2-C_1-C_4$ -alkylene-hetaryl;

R_A^3 :

each optionally substituted C_1-C_6 -alkyl, C_2-C_6 -alkenyl, C_2-C_6 -alkynyl, C_1-C_4 -alkylene- C_3-C_7 -cycloalkyl, C_1-C_4 -alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C_1-C_4 -alkylene-aryl, C_1-C_4 -alkylene-hetaryl, C_1-C_6 -alkylene-O- C_1-C_6 -alkyl, $CO-C_1-C_6$ -alkyl, CO -aryl, CO -hetaryl, $CO-C_1-C_4$ -alkylene-aryl, $CO-C_1-C_4$ -alkylene-hetaryl, $CO-O-C_1-C_6$ -alkyl, CO -aryl, $CO-O-C_1-C_4$ -alkylene-aryl, $CO-O$ -hetaryl, $CO-O-C_1-C_4$ -alkylene-

hetaryl, $\text{SO}_2\text{-C}_1\text{-C}_6\text{-alkyl}$, $\text{SO}_2\text{-aryl}$, $\text{SO}_2\text{-hetaryl}$, $\text{SO}_2\text{-C}_1\text{-C}_4\text{-alkylene-aryl}$ or $\text{SO}_2\text{-C}_1\text{-C}_4\text{-alkylene-hetaryl}$;

or the moieties $\mathbf{R_A}^2$ and $\mathbf{R_A}^3$ form, together with the nitrogen, a 3 to 7-membered, optionally substituted, saturated or aromatic heterocycle, which can contain one, two or three further different or identical heteroatoms from the group O, N, S; wherein optionally two moieties substituted on this heterocycle together can form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S, and wherein the so-formed cycle can optionally be substituted or a further, optionally substituted cycle can be condensed onto this cycle;

$\mathbf{R_A}^4$:

hydrogen, or
each optionally substituted $\text{C}_1\text{-C}_6\text{-alkyl}$, $\text{C}_1\text{-C}_6\text{-alkylene-O-C}_1\text{-C}_6\text{-alkyl}$, $\text{C}_2\text{-C}_6\text{-alkenyl}$, $\text{C}_3\text{-C}_{12}\text{-alkynyl}$, $\text{CO-C}_1\text{-C}_6\text{-alkyl}$, $\text{CO-O-C}_1\text{-C}_6\text{-alkyl}$, $\text{SO}_2\text{-C}_1\text{-C}_6\text{-alkyl}$, $\text{C}_3\text{-C}_7\text{-cycloalkyl}$, aryl, $\text{C}_1\text{-C}_4\text{-alkylene-aryl}$, CO-O-arylalkyl , $\text{CO-C}_1\text{-C}_4\text{-alkylene-aryl}$, CO-aryl , $\text{SO}_2\text{-aryl}$, hetaryl, CO-hetaryl or $\text{SO}_2\text{-C}_1\text{-C}_4\text{-alkylene-aryl}$;

B:

hydrogen or as moiety A is defined,

or each independently from another, two of the moieties A, B or $\mathbf{R_w}^1$ together form a 3 to 7-membered, optionally substituted, saturated or unsaturated carbocycle or an optionally substituted, saturated or unsaturated or aromatic heterocycle, which can contain one, two or three further different or identical heteroatoms from the group O, N, S; wherein optionally two moieties substituted on this carbo- or heterocycle can together form an anellated,

saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S and wherein the heterocycle formed can optionally be substituted or a further, optionally substituted cycle can be condensed onto this cycle;

R_w¹:

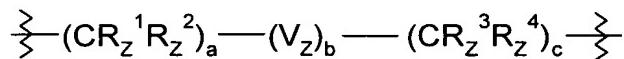
hydrogen, OH, halogen, NO₂, NH₂, CN, CF₃, CHF₂, O-CF₃, O-CHF₂, or each optionally substituted C₁-C₆-alkyl, C₃-C₇-cycloalkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₁-C₆-alkylene-S-C₁-C₆-alkyl, aryl, hetaryl, O-C₁-C₆-alkyl, O-aryl, O-benzyl, C₁-C₆-alkylamino, C₁-C₆-dialkylamino, pyrrolidinyl, piperidinyl, morpholinyl, CO-C₁-C₆-alkyl, SO₂-C₁-C₆-alkyl, CO-aryl, SO₂-aryl, CO-C₁-C₄-alkylene-aryl, SO₂-C₁-C₄-alkylene-aryl, SO-aryl, CONH₂, CONH-C₁-C₆-alkyl, SO₂NH-C₁-C₆-alkyl, CON(C₁-C₆-alkyl)₂, SO₂N(C₁-C₆-alkyl)₂, NH-SO₂-C₁-C₆-alkyl or NH-CO-C₁-C₆-alkyl;

D:

as moiety A is defined;

Z:

a moiety of the general formula **Z1**



Z1

with the indices

$$a = 0 - 4$$

$$b = 0, 1$$

$c = 0 - 4$

wherein the sum of a, b and c is no more than 5;

$R_Z^1, R_Z^2, R_Z^3, R_Z^4$ independently of one another:

hydrogen, halogen, OH, or

each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₆-alkylene-C₃-C₇-cycloalkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, hetaryl or C₁-C₄-alkylene-hetaryl, or

each independently of one another, two moieties R_Z^1 and R_Z^2 or R_Z^3 and R_Z^4 together form a 3 to 7-membered, optionally substituted, saturated or unsaturated carbo- or heterocycle, which can contain up to three heteroatoms from the group O, N or S;

V_Z :

-CO-, -CO-NR_z⁵-, -NR_z⁵-CO-, -O-, -S-, -SO-, -SO₂-, -SO₂-NR_z⁵-, -NR_z⁵-SO₂-, -CS-, -CS-NR_z⁵-, -NR_z⁵-CS-, -CS-O-, -O-CS-, -CO-O-, -O-CO-, -O-, ethynylene, -C(=CR_z⁶R_z⁷)-, -CR_z⁶=CR_z⁷-, -NR_z⁵-CO-NR_z^{5*}-, -O-CO-NR_z⁵-, -NR_z⁵-;

R_Z^5, R_Z^{5*} independently of one another:

hydrogen, or

each optionally substituted C₁-C₆-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₂-C₆-alkenyl, C₃-C₁₂-alkynyl, CO-C₁-C₆-alkyl, CO-O-C₁-C₆-alkyl, SO₂-C₁-C₆-alkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-aryl, CO-aryl, SO₂-aryl, hetaryl, CO-hetaryl or SO₂-C₁-C₄-alkylene-aryl;

R_Z⁶, R_Z⁷ independently of one another:

hydrogen, OH, or

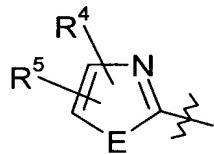
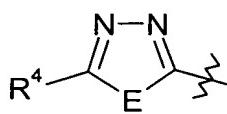
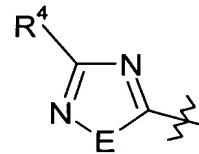
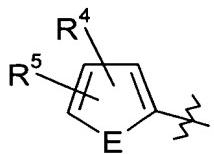
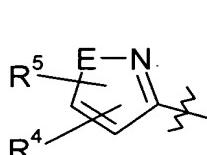
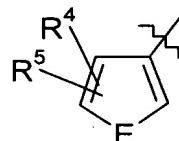
each optionally substituted C₁-C₆-alkyl, C₁-C₄-alkoxy, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₆-alkylene-C₃-C₇-cycloalkyl, C₃-C₇-cycloalkyl, aryl, C₁-C₄-alkylene-aryl, hetaryl or C₁-C₄-alkylene-hetaryl;

R¹, R², R³ independently of one another:

hydrogen, OH, CN, or

each optionally substituted C₁-C₆-alkyl, O-C₁-C₆-alkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₃-C₇-cycloalkyl, O-C₃-C₇-cycloalkyl, aryl, hetaryl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, O-aryl, O-C₁-C₄-alkylene-aryl, O-hetaryl, O-C₁-C₄-alkylene-hetaryl, CO-C₁-C₆-alkyl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-aryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl, OCO-C₁-C₆-alkyl, OCO-aryl, OCO-hetaryl, OCO-C₁-C₄-alkylene-aryl, OCO-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl or SO₂-C₁-C₄-alkylene-aryl, or

each independently of the third moiety, two moieties of R¹, R² or R³ together form a 5 to 7-membered, optionally substituted, saturated or unsaturated carbocycle, or an optionally substituted, saturated or unsaturated, which can contain one, two or three different or identical heteroatoms from the group O, N, S, wherein optionally two moieties substituted on this carbo- or heterocycle can together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle may contain up to three different or identical heteroatoms O, N, S and wherein the cycle formed can optionally be substituted or a further, optionally substituted cycle can be condensed onto this cycle;

Q:a doubly substituted 5-membered hetaryl moiety, chosen from **Q1** to **Q6****Q1****Q2****Q3****Q4****Q5****Q6**E: O, N-R_Q¹ or S;R_Q¹:

hydrogen, or

each optionally substituted C₁-C₄-alkyl, CO-C₁-C₄-alkyl, SO₂-C₁-C₄-alkyl,
 CO-O-C₁-C₄-alkyl, aryl, C₁-C₄-alkylene-aryl, CO-aryl, CO-hetaryl, SO₂-
 aryl, SO₂-hetaryl, CO-O-aryl, CO-C₁-C₄-alkylene-aryl, SO₂-C₁-C₄-
 alkylene-aryl or CO-O-C₁-C₄-alkylene-aryl;

R⁴, R⁵ each independently of one another a moiety chosen from the groups 1.),
 2.), 3.), 4.) or 5.):

- 1.) hydrogen, halogen, CN, CF₃, CHF₂, or
 each optionally substituted C₁-C₁₀-alkyl, C₂-C₁₀-alkenyl, C₂-C₁₀-alkynyl, C₃-C₇-cycloalkyl, C₁-C₆-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, C₁-C₆-alkylene-O-aryl, COO-C₁-C₄-alkyl or C₁-C₄-alkylene-COO-C₁-C₄-alkyl;

[[4]] 2.) Phenyl or naphthyl, which are each substituted with R_Q², R_Q³ and R_Q⁴,

wherein

R_Q², R_Q³ and R_Q⁴ each independently from one another represent a substituent from the following group:

hydrogen, NO₂, NH₂, OH, CN, CF₃, CHF₂, OCF₃, OCHF₂, COOH, O-CH₂-COOH, SH, halogen, or
 each optionally substituted aryl, hetaryl, heterocycloalkyl, C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, C₁-C₄-alkylene-aryl or C₁-C₄-alkylene-hetaryl, or
 O-R_Q⁵, S-R_Q⁵, NR_Q⁷R_Q⁸, CO-OR_Q⁶, NR_Q⁸-CO-O-R_Q⁶, O-CH₂-COO-R_Q⁶, NR_Q⁸-CO-R_Q⁶, SO₂-R_Q⁶, NR_Q⁸-SO₂-R_Q⁶, SO₂NH₂, CONH₂, SO₂-NR_Q⁷R_Q⁸ or CO-NR_Q⁷R_Q⁸, or

two of the moieties R_Q², R_Q³ or R_Q⁴ together form a 3 to 7-membered, optionally substituted, saturated, unsaturated carbocycle or an optionally substituted, saturated, unsaturated aromatic heterocycle, which can contain up to three further different or identical heteroatoms O, N, S, and optionally two moieties substituted on this heterocycle together can form an anellated, saturated, unsaturated or aromatic

carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S, and the cycle formed can optionally be substituted or a further, optionally substituted cycle can be condensed onto this cycle;

R_Q⁵ each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, heterocycloalkyl, aryl or hetaryl;

R_Q⁶ each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl or C₁-C₆-alkylene-O-C₁-C₆-alkyl;

R_Q⁷ hydrogen, OH, CN, or
each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, CO-C₁-C₆-alkyl, C₁-C₄-alkylene-aryl, C₁-C₄-alkylene-hetaryl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-O-hetaryl, CO-O-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl or SO₂-C₁-C₄-alkylene-hetaryl;

R_Q⁸ each optionally substituted C₁-C₆-alkyl, C₂-C₆-alkenyl, C₂-C₆-alkynyl, C₃-C₇-cycloalkyl, C₁-C₄-alkylene-C₃-C₇-cycloalkyl, C₁-C₄-alkylene-heterocycloalkyl, aryl, hetaryl, heterocycloalkyl, C₁-C₆-alkylene-O-C₁-C₆-alkyl, CO-C₁-C₆-alkyl, CO-aryl, CO-hetaryl, CO-C₁-C₄-alkylene-aryl, CO-C₁-C₄-alkylene-hetaryl, CO-O-C₁-C₆-alkyl, CO-O-aryl, CO-O-C₁-C₄-alkylene-aryl, CO-O-

hetaryl, CO-O-C₁-C₄-alkylene-hetaryl, SO₂-C₁-C₆-alkyl, SO₂-aryl, SO₂-hetaryl, SO₂-C₁-C₄-alkylene-aryl or SO₂-C₁-C₄-alkylene-hetaryl;

or both moieties R_Q⁷ and R_Q⁸ form, together with the nitrogen, a 3 to 7-membered, optionally substituted, saturated or aromatic heterocycle, which can contain one, two or three further different or identical heteroatoms O, N, S; and optionally two moieties substituted on this heterocycle can together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or same heteroatoms O, N, S and the cycle formed can optionally be substituted or a further, optionally substituted cycle can be condensed onto this cycle;

- 3.) a 5- or 6-membered, hetaryl moiety, optionally substituted with one or two substituents from the group consisting of:

2-furyl, 3-furyl, 2-pyrrolyl, 3-pyrrolyl, 2-thienyl, 3-thienyl, 2-pyridyl, 3-pyridyl, 4-pyridyl, 2-thiazolyl, 4-thiazolyl, 5-thiazolyl, 2-oxazolyl, 4-oxazolyl, 5-oxazolyl, 2-pyrimidyl, 4-pyrimidyl, 5-pyrimidyl, 6-pyrimidyl, 3-pyrazolyl, 4-pyrazolyl, 5-pyrazolyl, 3-isothiazolyl, 4-isothiazolyl, 5-isothiazolyl, 2-imidazolyl, 4-imidazolyl, 5-imidazolyl, 3-pyridazinyl, 4-pyridazinyl, 5-pyridazinyl, 6-pyridazinyl, 3-isoxazolyl, 4-isoxazolyl, 5-isoxazolyl, thiadiazolyl, oxadiazolyl or triazinyl or their anellated derivatives indazolyl, indolyl, benzothiophenyl, benzofuranyl, indolinyl, benzimidazolyl, benzthiazolyl, benzoxazolyl, chinolinyl and isochinolinyl;

- 4.) both moieties R⁴ and R⁵ together form a 4 to 7-membered, optionally substituted, saturated or unsaturated or aromatic carbocycle or a 5- or 6-

membered optionally substituted, saturated or unsaturated or aromatic heterocycle, which can contain up to three further different or identical heteroatoms O, N, S, and can be substituted with up to two further moieties, wherein optionally two moieties substituted on this carbo- or heterocycle can together form an anellated, saturated, unsaturated or aromatic carbocycle or heterocycle, wherein the heterocycle can contain up to three different or identical heteroatoms O, N, S and wherein the cycle formed can optionally be substituted or a further, optionally substituted cycle can be condensed onto this cycle;

- 5.) a C₆-C₁₀- bi- or tricyclic, saturated hydrocarbon moiety;

~~for the manufacture of a medicament for the treatment of diseases which are modulated by a 5-HT5 receptor activity.~~

18. (Currently Amended) Use A method according to claim 16, wherein R⁴ and/or R⁵ have the following meanings:

2-pyridyl, 3-pyridyl, 4-pyridyl, 2-thienyl, 3-thienyl, benzothiophenyl, benzofuranyl, chinolinyl or isochinolinyl, which may optionally be substituted with 1 or 2 moieties.

19. (Currently Amended) Use A method according to claim 16 or 17 ~~for the treatment of where the disease is characterized by neuropathological, neuropsychiatric and neurodegenerative disorders, symptoms and dysfunctions.~~

20. (Currently Amended) Use A method according to at least one of the claims claim 16 to 18 for the treatment of where the disease is characterized by migraine and brain damage.
21. (Currently Amended) Use A method according to claim 18 for the treatment of neuropathological, neuropsychiatric and neurodegenerative diseases, selected from the group consisting of cerebral ischemia, stroke, epilepsy and seizures in general, psychoses, schizophrenia, autism, OCD-syndrome, cognitive diseases, attention disorders, depressions, bipolar- and/or unipolar depressions, states of anxiety, dementia, senile dementia, Alzheimer dementia, demyelinizing diseases, multiple sclerosis and brain tumors.
22. (Currently Amended) Use A method according to claim 16 or 17 for the treatment of diseases chosen from the group consisting of cerebral vascular disorders, pain, disorders due to pain, addiction, disorders due to drugs, amnesia, alcohol abuse, drug abuse, disorders of the circadian rhythm and Cushing syndrome.